

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM042822\  
 Data File : BM034865.D  
 Acq On : 28 Apr 2022 14:02  
 Operator : CG/JU  
 Sample : PB144441BL  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 PB144441BL

Quant Time: Apr 29 05:06:01 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM042722.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Apr 28 01:51:59 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.928  | 152  | 213412   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.734 | 136  | 812230   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.563 | 164  | 448129   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.304 | 188  | 815440   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.480 | 240  | 597218   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 23.868 | 264  | 548606   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.505  | 112  | 1750170  | 144.682 | ng    | 0.00     |
| 7) Phenol-d6                | 7.093  | 99   | 2113861  | 135.364 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.087  | 82   | 1401816  | 93.201  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.045 | 330  | 608328   | 124.792 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.192 | 172  | 2994113  | 96.992  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.927 | 244  | 3352090  | 108.676 | ng    | 0.00     |

|                  |        |
|------------------|--------|
| Target Compounds | Qvalue |
| -----            |        |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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